

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
 Data File : BN019637.D
 Acq On : 03 May 2022 01:43
 Operator : CG/JU
 Sample : N2543-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ETNS5

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
 Title : SVOA CALIBRATION

Signal : TIC: BN019637.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.187	14	17	20	rBV2	59473	66660	1.59%	0.399%
2	3.217	20	22	24	rVV2	27270	29672	0.71%	0.178%
3	3.240	24	26	28	rVV2	43930	42977	1.03%	0.257%
4	3.281	28	33	43	rVB	3524306	4190980	100.00%	25.109%
5	3.581	81	84	85	rBV2	7445	7071	0.17%	0.042%
6	3.611	85	89	96	rVB	98392	138356	3.30%	0.829%
7	3.705	101	105	110	rVB3	10137	16249	0.39%	0.097%
8	4.081	165	169	174	rVB	13493	17460	0.42%	0.105%
9	5.199	354	359	364	rVB5	3385	6250	0.15%	0.037%
10	7.869	806	813	822	rBV	584287	942720	22.49%	5.648%
11	10.457	1245	1253	1258	rBV7	2933	6082	0.15%	0.036%
12	10.657	1277	1287	1298	rBV	712999	1269988	30.30%	7.609%
13	10.804	1308	1312	1315	rVB4	3735	5286	0.13%	0.032%
14	11.140	1363	1369	1376	rVB8	3711	7401	0.18%	0.044%
15	13.339	1737	1743	1745	rBV4	3479	4297	0.10%	0.026%
16	14.486	1931	1938	1946	rBV	954602	1430773	34.14%	8.572%
17	14.651	1959	1966	1970	rBV5	5138	8540	0.20%	0.051%
18	15.739	2146	2151	2155	rBV4	11450	16423	0.39%	0.098%
19	16.763	2319	2325	2330	rBV2	38153	54513	1.30%	0.327%
20	17.010	2364	2367	2370	rBV4	4353	5688	0.14%	0.034%
21	17.110	2379	2384	2388	rBV	236340	313844	7.49%	1.880%
22	17.145	2388	2390	2395	rVB	50785	60226	1.44%	0.361%
23	17.228	2397	2404	2415	rBV	1027017	1544636	36.86%	9.254%
24	17.404	2429	2434	2441	rBV	148810	227647	5.43%	1.364%
25	17.686	2477	2482	2486	rBV2	29653	44056	1.05%	0.264%
26	17.833	2499	2507	2514	rBV	293010	507763	12.12%	3.042%
27	17.916	2518	2521	2523	rBV4	3710	5546	0.13%	0.033%
28	17.951	2523	2527	2533	rVB4	32265	50332	1.20%	0.302%
29	18.027	2535	2540	2544	rBV7	7901	10929	0.26%	0.065%
30	18.080	2544	2549	2554	rBV	113348	150224	3.58%	0.900%
31	18.139	2554	2559	2563	rVB3	30372	39907	0.95%	0.239%
32	18.239	2572	2576	2582	rBV6	11417	17971	0.43%	0.108%
33	18.304	2582	2587	2591	rVV	93126	127618	3.05%	0.765%
34	18.363	2593	2597	2601	rVV2	70864	91219	2.18%	0.547%
35	18.410	2601	2605	2611	rVB2	53411	74722	1.78%	0.448%
36	18.586	2630	2635	2640	rBV	99389	145813	3.48%	0.874%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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37	18.633	2640	2643	2647	rVV3	37842	50424	1.20%	0.302%
38	18.686	2647	2652	2656	rVV	66489	90965	2.17%	0.545%
39	18.727	2656	2659	2662	rVV4	32105	45376	1.08%	0.272%
40	18.763	2662	2665	2669	rVB2	74967	93048	2.22%	0.557%
41	18.863	2678	2682	2686	rVB2	27724	33676	0.80%	0.202%
42	19.069	2713	2717	2722	rBV	47312	59414	1.42%	0.356%
43	19.122	2722	2726	2729	rVV2	88342	113961	2.72%	0.683%
44	19.157	2729	2732	2737	rVB	88416	121534	2.90%	0.728%
45	19.269	2749	2751	2755	rVB5	6125	5964	0.14%	0.036%
46	19.374	2763	2769	2775	rBV4	69089	150358	3.59%	0.901%
47	19.433	2775	2779	2782	rVB4	13776	18354	0.44%	0.110%
48	19.492	2787	2789	2793	rBV5	6305	6921	0.17%	0.041%
49	19.533	2793	2796	2799	rBV4	6162	7228	0.17%	0.043%
50	19.763	2830	2835	2840	rBV	169980	213890	5.10%	1.281%
51	19.863	2848	2852	2859	rBV	122449	162814	3.88%	0.975%
52	19.922	2859	2862	2866	rVV5	9282	10295	0.25%	0.062%
53	20.269	2916	2921	2926	rBV	232206	281078	6.71%	1.684%
54	20.321	2926	2930	2935	rVV	112295	146735	3.50%	0.879%
55	20.710	2991	2996	3001	rBV	852013	983388	23.46%	5.892%
56	21.410	3110	3115	3120	rBV	932586	1228381	29.31%	7.359%
57	23.757	3508	3514	3528	rVB2	581064	1187775	28.34%	7.116%

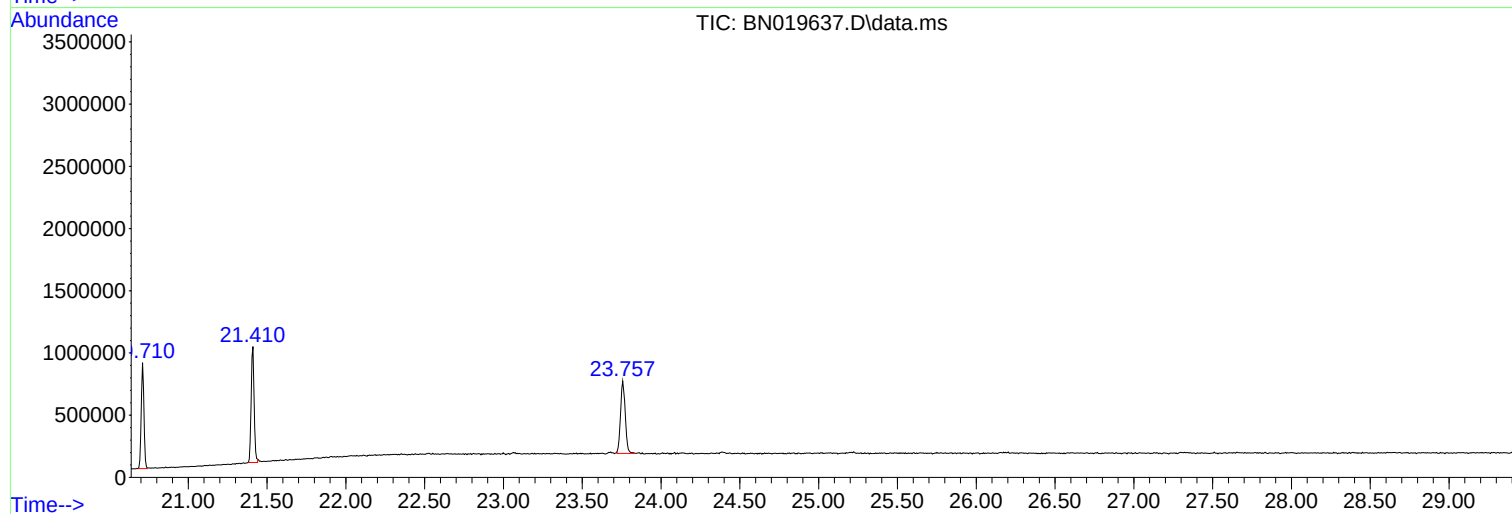
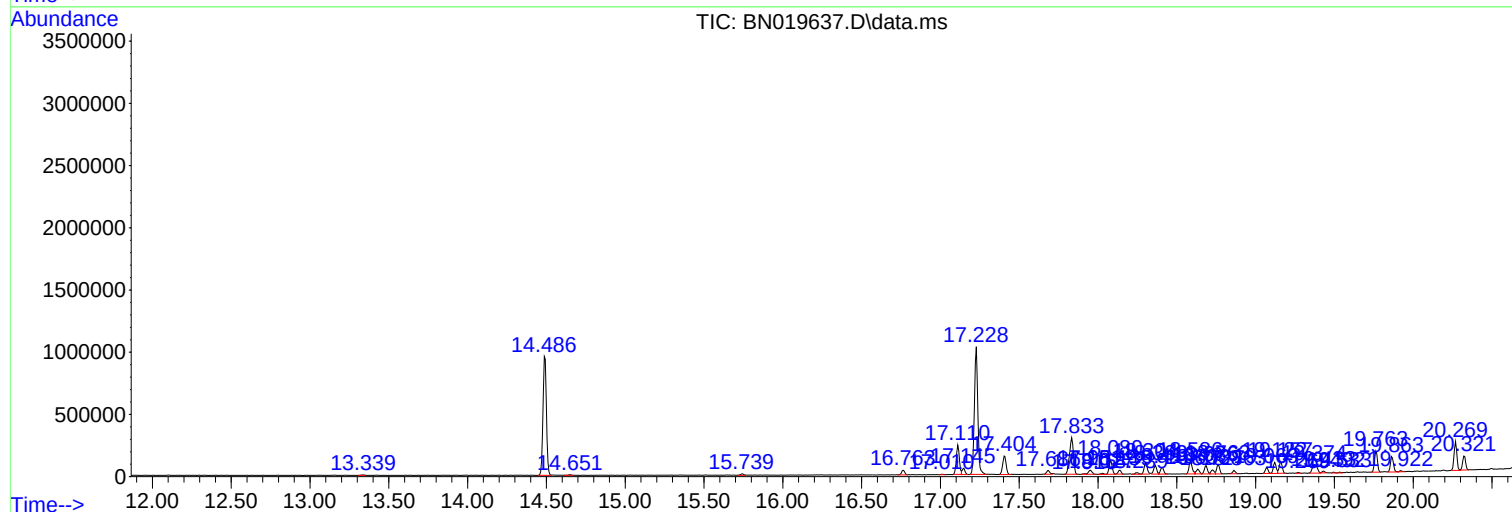
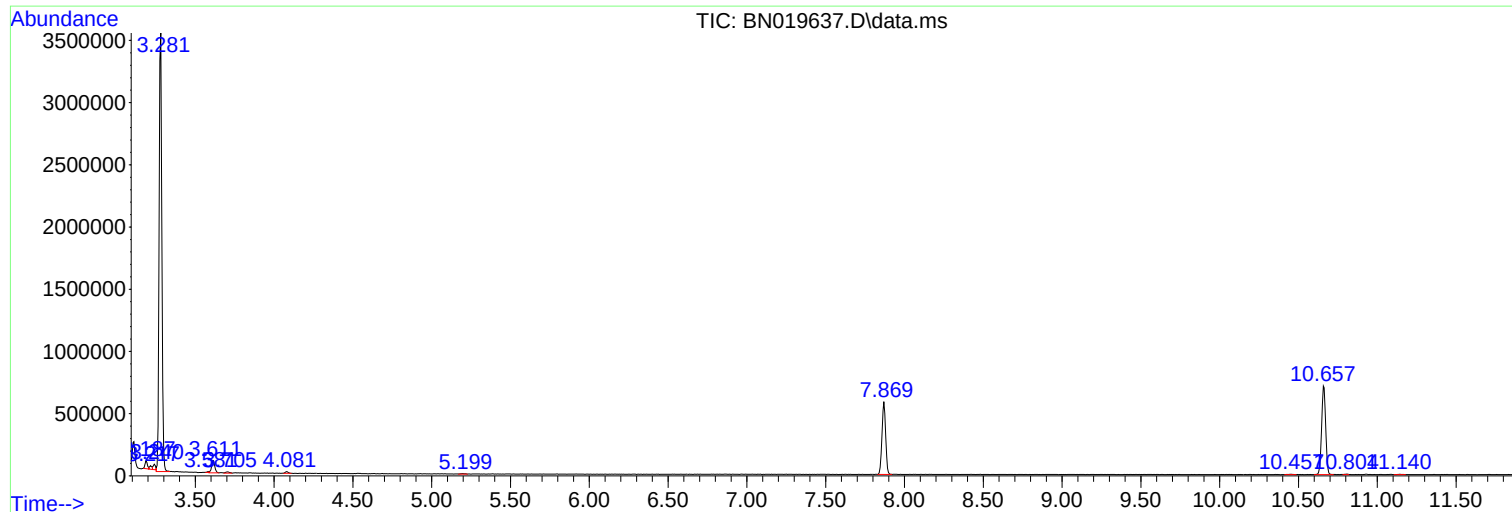
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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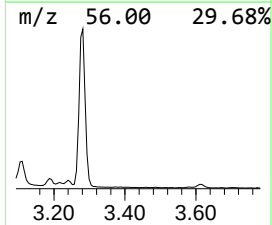
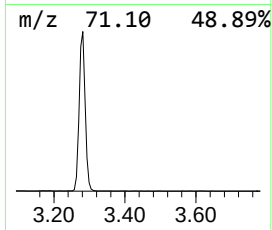
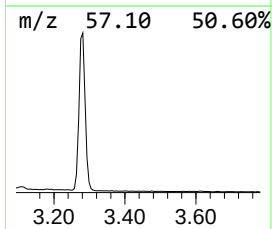
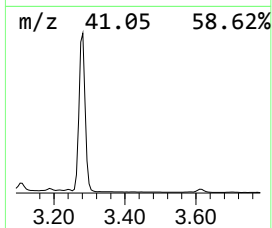
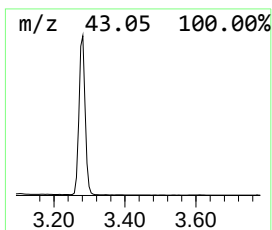
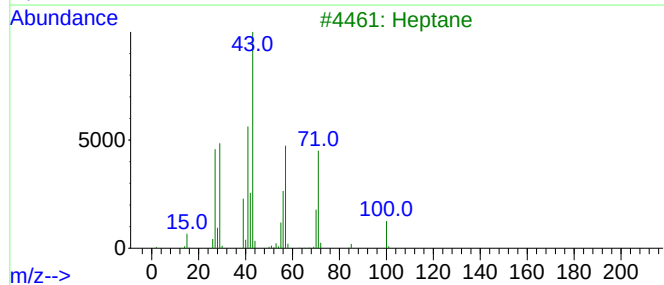
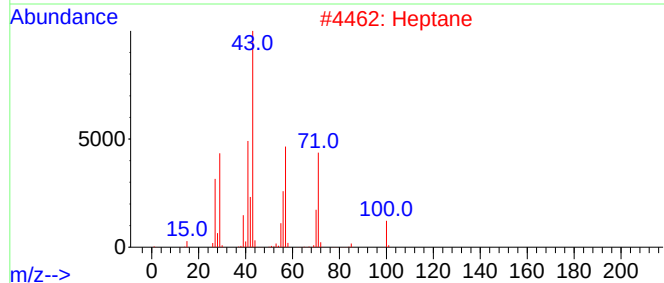
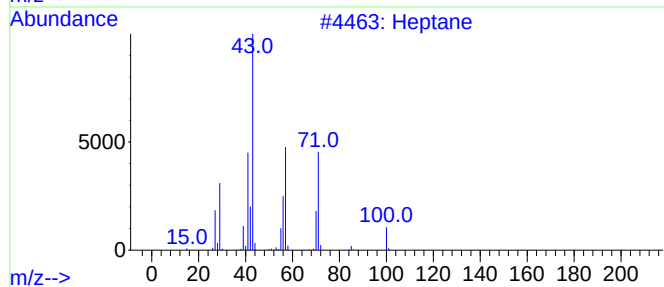
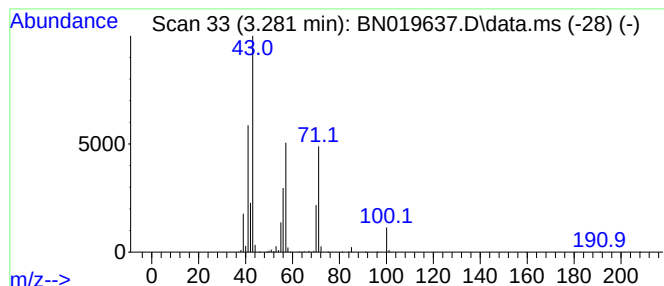
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.281	88.91 ng/ul	4190980	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane			100	C7H16	000142-82-5	95
2	Heptane			100	C7H16	000142-82-5	94
3	Heptane			100	C7H16	000142-82-5	91
4	Heptane			100	C7H16	000142-82-5	83
5	Oxalic acid, isobutyl pentyl ester			216	C11H20O4	1000309-37-0	64



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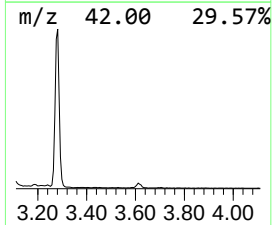
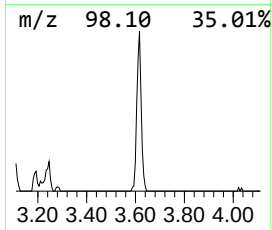
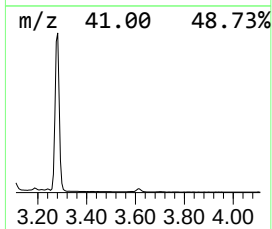
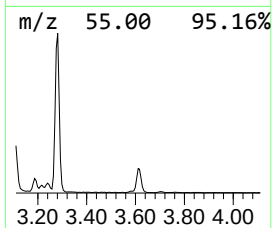
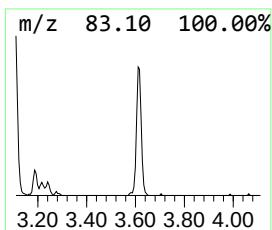
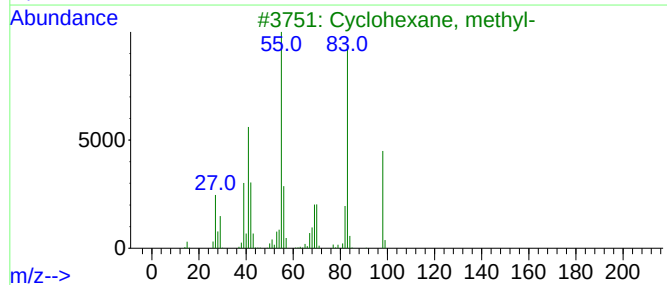
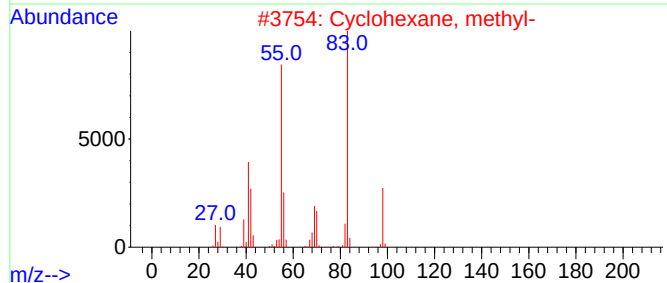
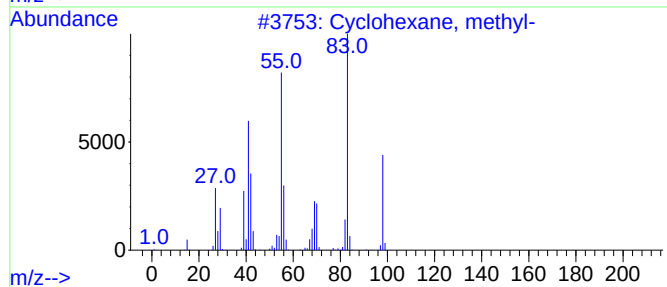
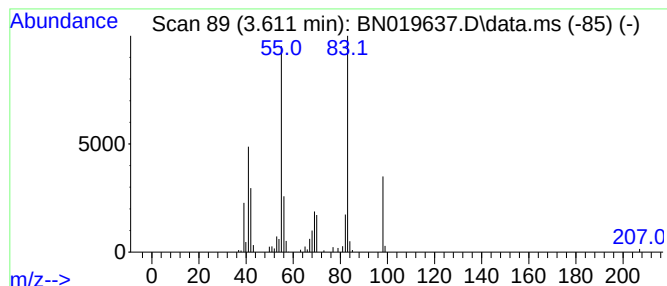
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Cyclic3.611 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.611	2.94 ng/ul	138356	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, methyl-	98	C7H14	000108-87-2	94
2			Cyclohexane, methyl-	98	C7H14	000108-87-2	94
3			Cyclohexane, methyl-	98	C7H14	000108-87-2	94
4			Cyclohexane, methyl-	98	C7H14	000108-87-2	91
5			Cyclohexane, methyl-	98	C7H14	000108-87-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
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ALS Vial : 26 Sample Multiplier: 1

Instrument :
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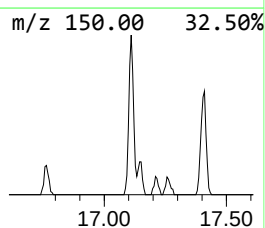
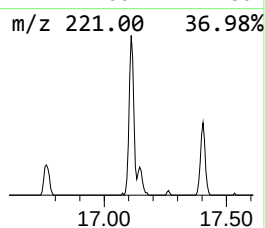
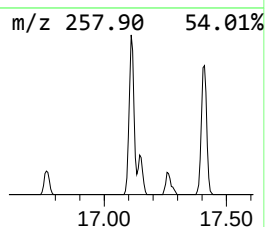
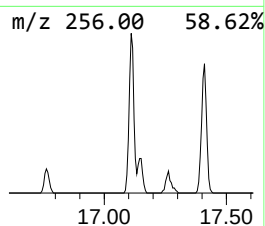
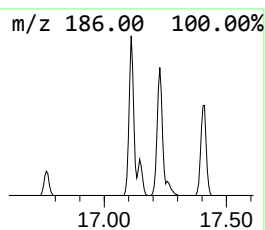
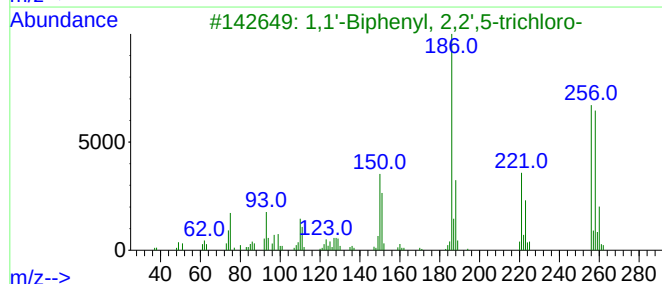
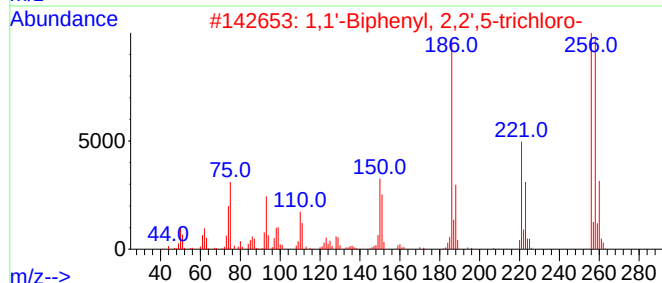
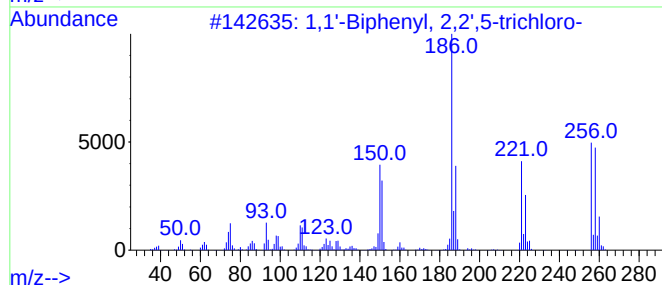
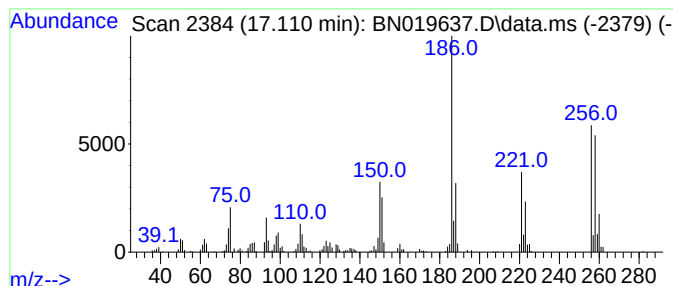
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1,1'-Biphenyl, 2,2',5-trich... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.110	4.06 ng/ul	313844	Phenanthrene-d10	17.227

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
2	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
3	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
4	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
5	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		



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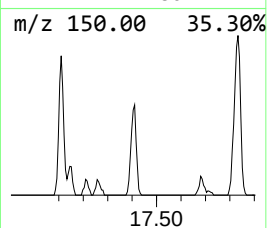
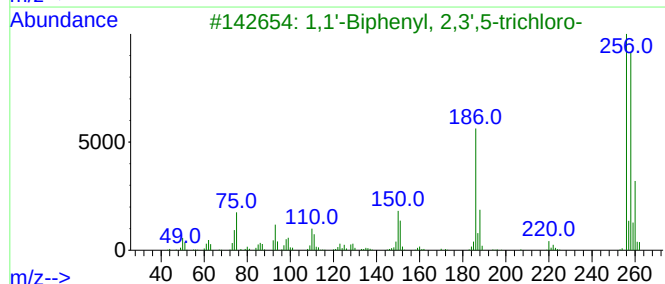
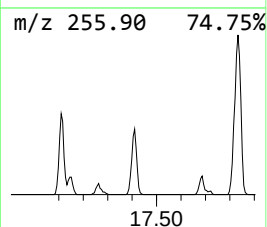
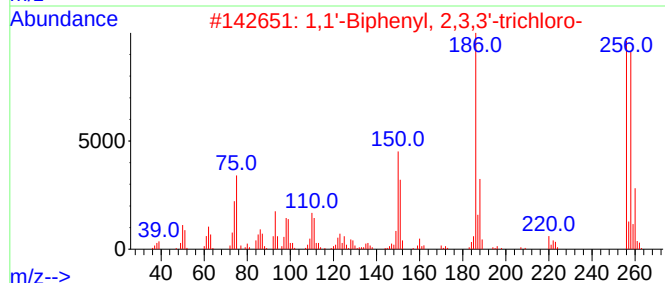
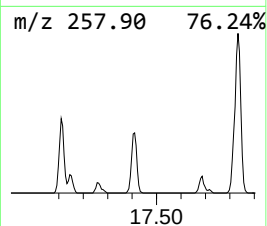
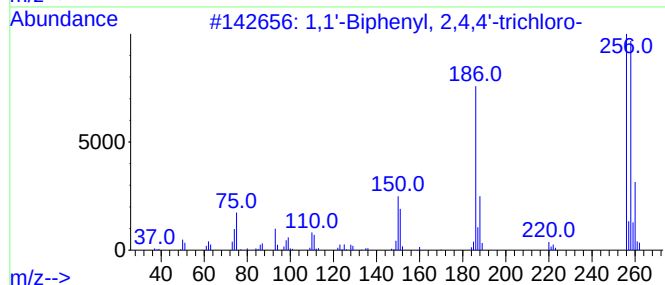
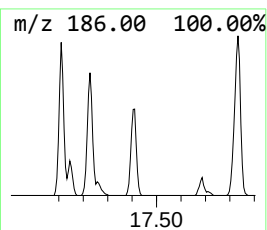
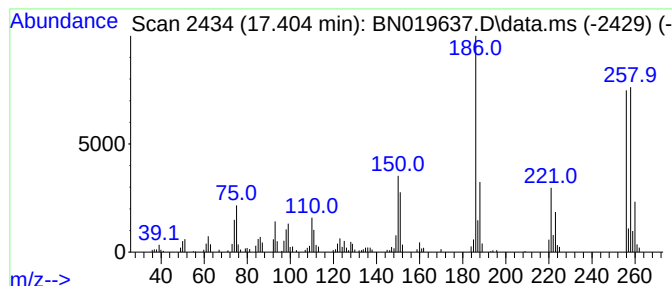
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.404	2.95 ng/ul	227647	Phenanthrene-d10	17.227

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
3	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99		
4	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99		
5	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99		



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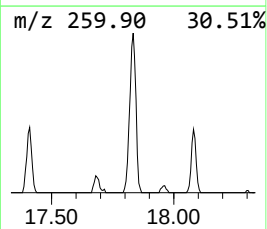
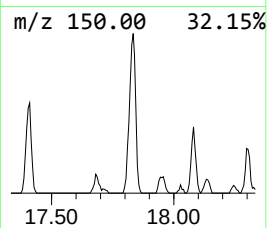
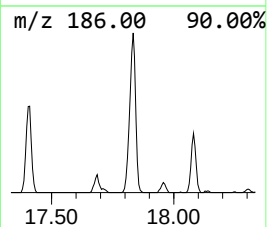
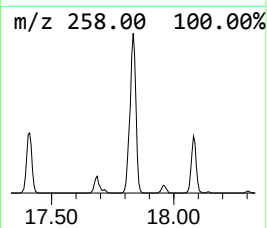
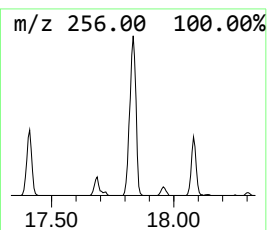
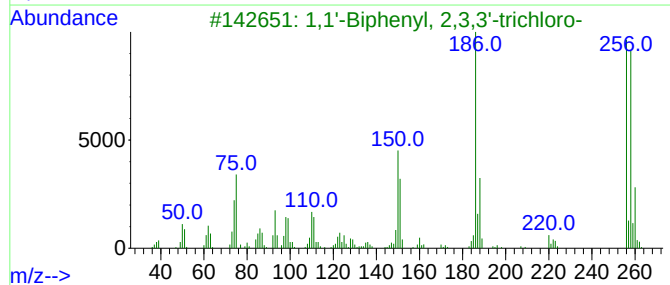
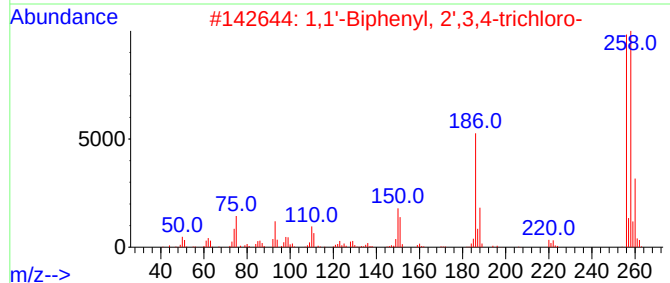
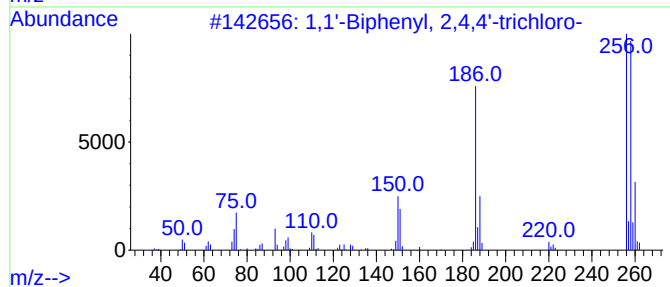
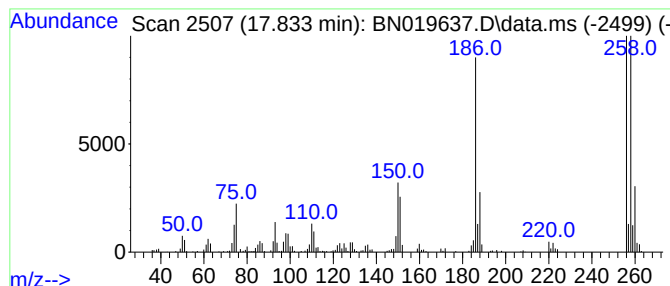
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1,1'-Biphenyl, 2',3,4-trich... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.833	6.57 ng/ul	507763	Phenanthrene-d10	17.227

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
2	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99		
3	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
4	1,1'-Biphenyl, 2',3,5-trichloro-	256	C12H7Cl3	037680-68-5	99		
5	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
Data File : BN019637.D
Acq On : 03 May 2022 01:43
Operator : CG/JU
Sample : N2543-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

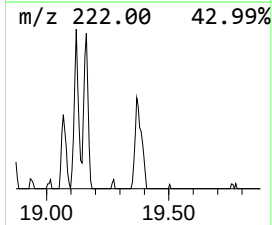
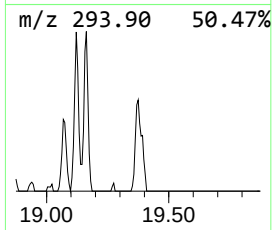
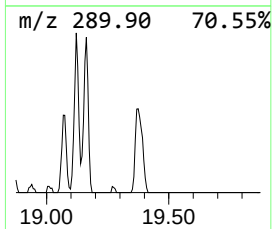
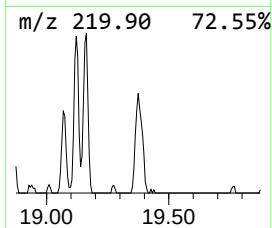
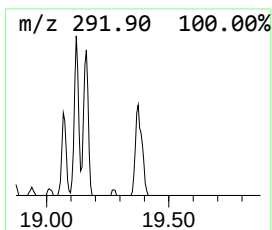
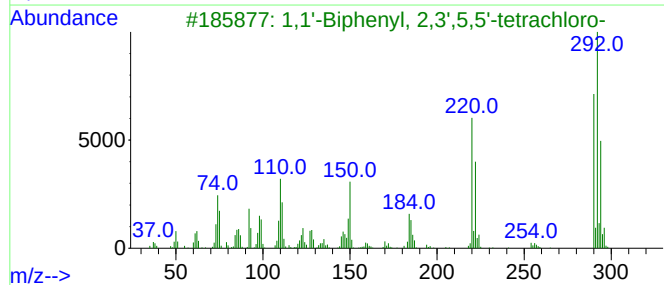
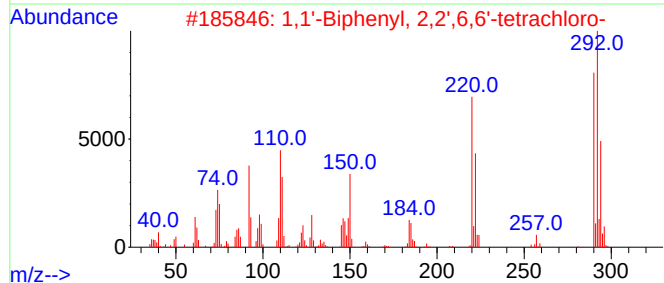
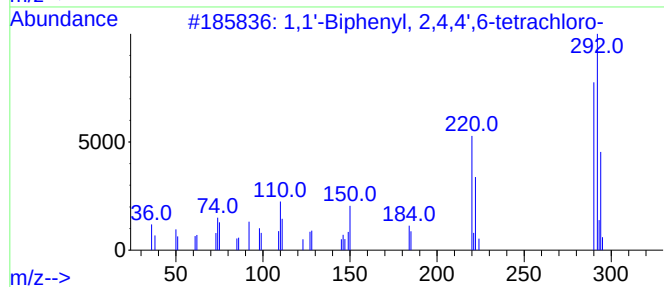
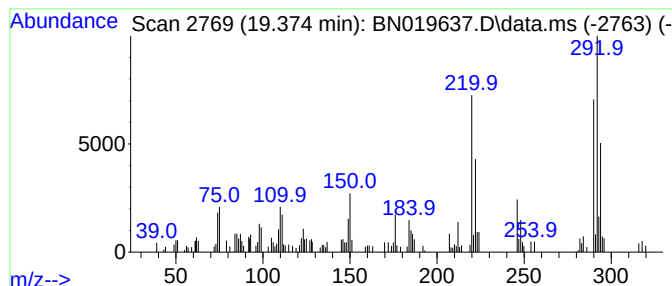
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1,1'-Biphenyl, 2,4,4',6-tet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.375	2.45 ng/ul	150358	Chrysene-d12	21.410	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4',6-tetrachl...	290	C12H6Cl4	032598-12-2	99
2	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	98
3	1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	98
4	1,1'-Biphenyl, 3,3',4,5'-tetrach...	290	C12H6Cl4	041464-48-6	97
5	1,1'-Biphenyl, 2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
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Acq On : 03 May 2022 01:43
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Instrument :
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ClientSampleId :
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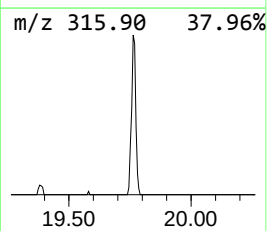
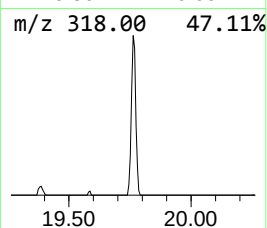
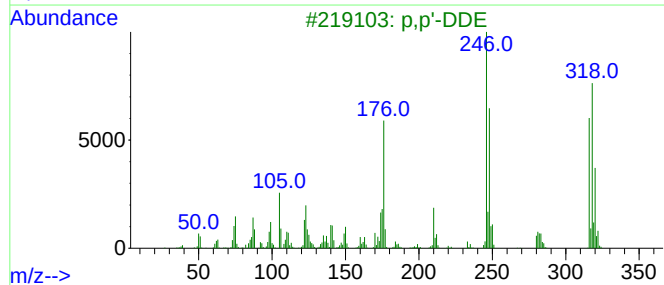
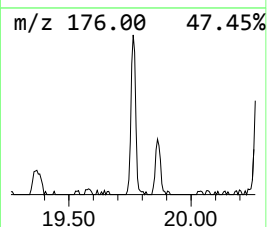
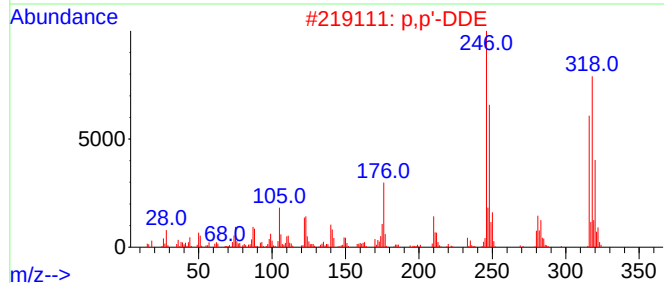
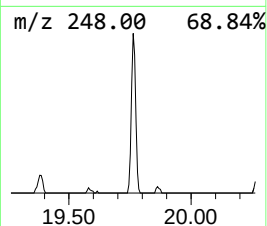
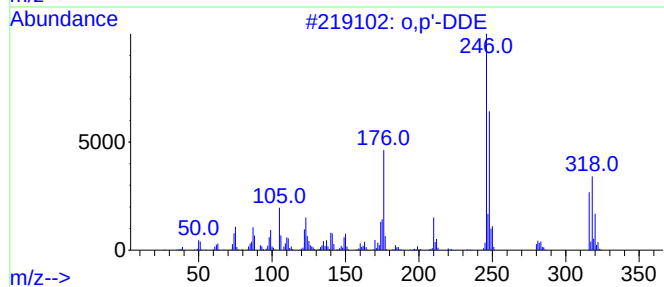
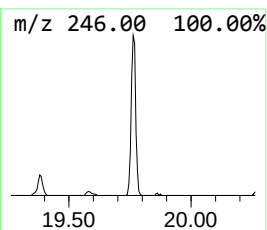
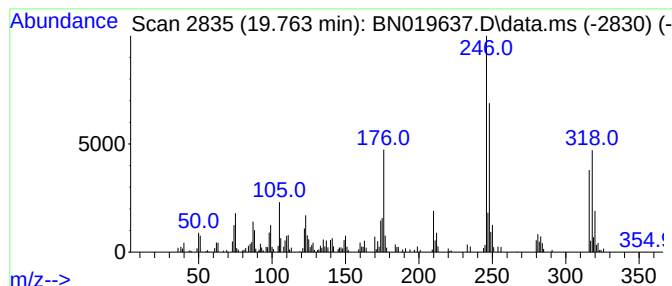
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 o,p'-DDE Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.763	3.48 ng/ul	213890	Chrysene-d12	21.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o,p'-DDE	316	C14H8Cl4	003424-82-6	99
2			p,p'-DDE	316	C14H8Cl4	000072-55-9	99
3			p,p'-DDE	316	C14H8Cl4	000072-55-9	99
4			p,p'-DDE	316	C14H8Cl4	000072-55-9	99
5			p,p'-DDE	316	C14H8Cl4	000072-55-9	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
Data File : BN019637.D
Acq On : 03 May 2022 01:43
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Sample : N2543-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

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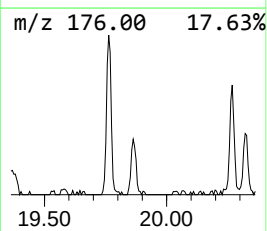
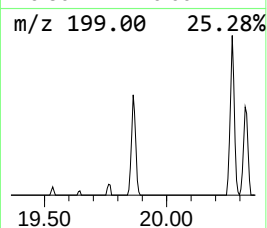
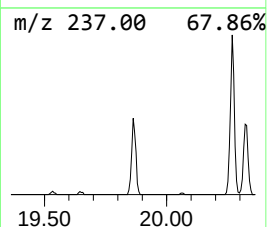
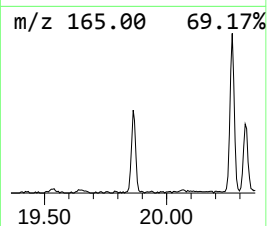
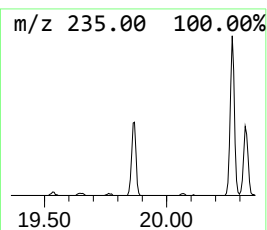
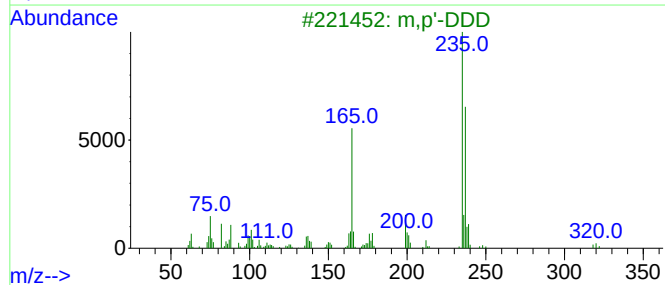
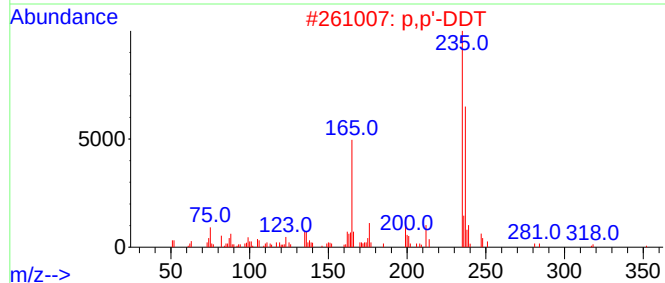
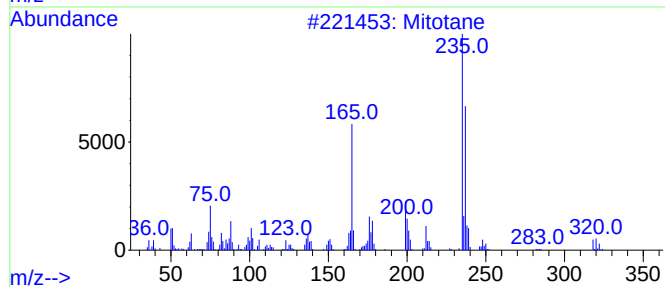
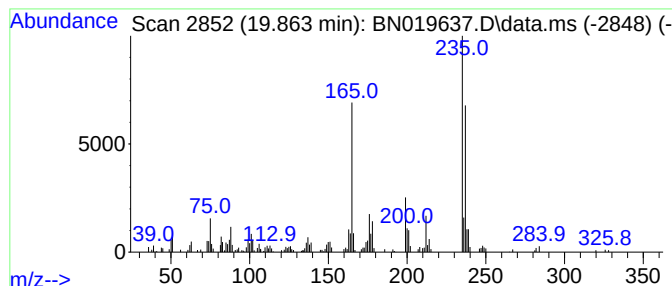
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Mitotane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.863	2.65 ng/ul	162814	Chrysene-d12	21.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mitotane	318	C14H10Cl4	000053-19-0	93
2			p,p'-DDT	352	C14H9Cl5	000050-29-3	91
3			m,p'-DDD	318	C14H10Cl4	004329-12-8	90
4			Mitotane	318	C14H10Cl4	000053-19-0	90
5			1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
Data File : BN019637.D
Acq On : 03 May 2022 01:43
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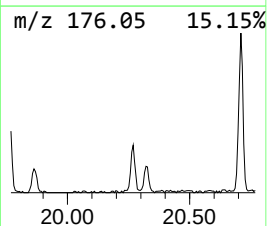
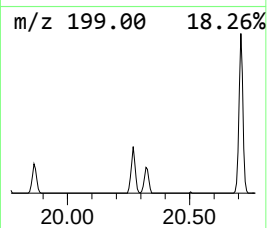
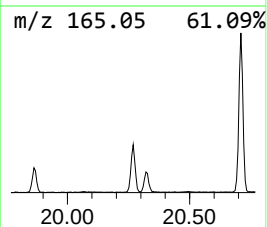
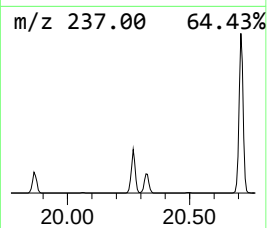
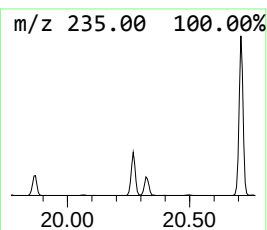
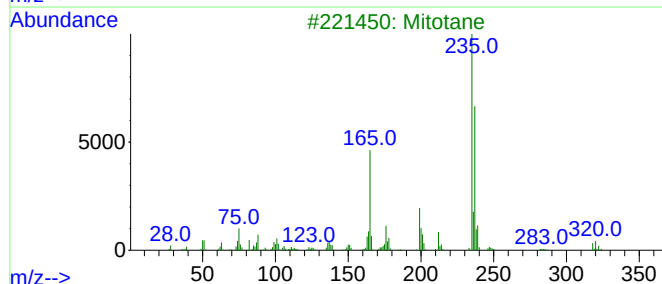
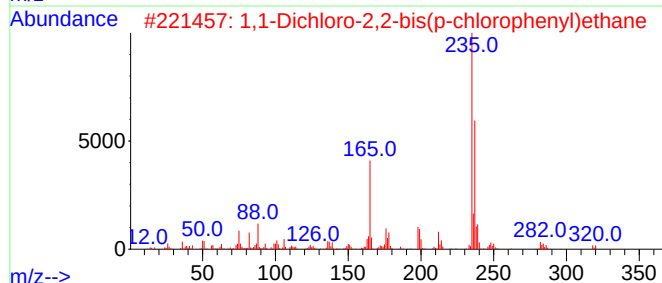
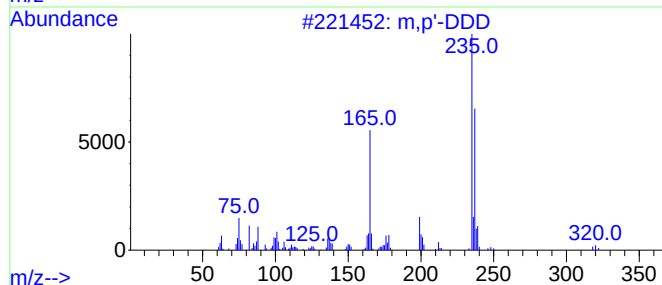
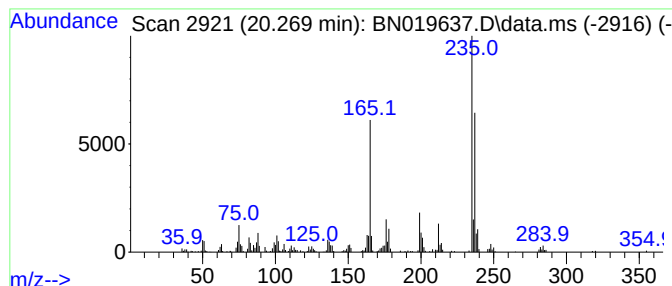
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 m,p'-DDD Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.269	4.58 ng/ul	281078	Chrysene-d12	21.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	m,p'-DDD			318	C14H10Cl4	004329-12-8	98
2	1,1-Dichloro-2,2-bis(p-chlorophe...			318	C14H10Cl4	000072-54-8	97
3	Mitotane			318	C14H10Cl4	000053-19-0	91
4	Mitotane			318	C14H10Cl4	000053-19-0	91
5	1,1-Dichloro-2,2-bis(p-chlorophe...			318	C14H10Cl4	000072-54-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
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Sample : N2543-01
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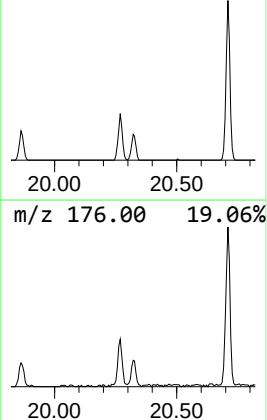
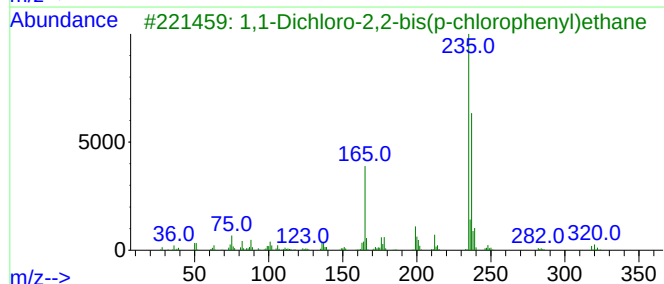
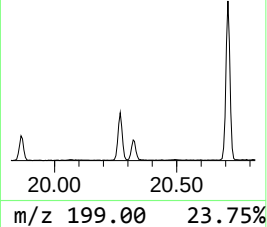
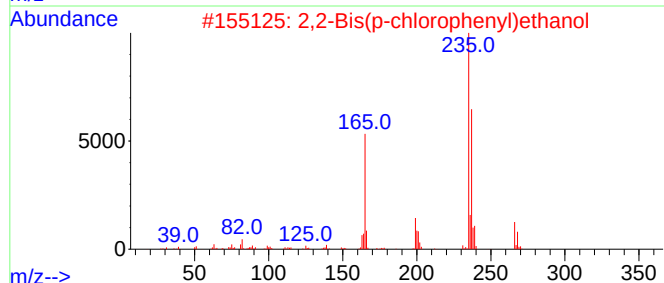
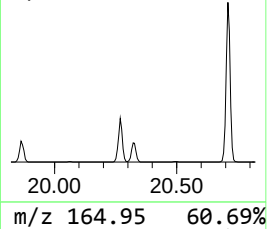
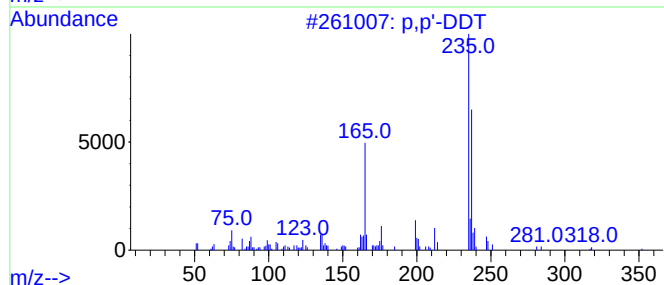
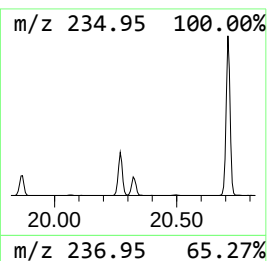
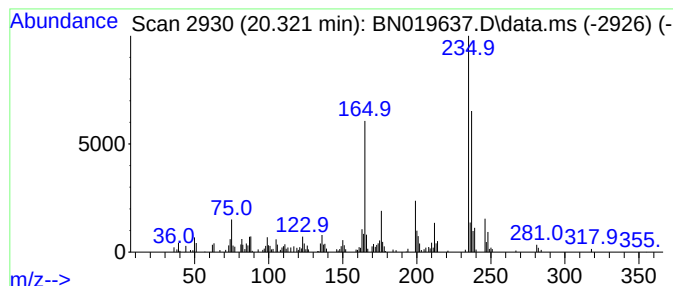
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 p,p'-DDT Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.322	2.39 ng/ul	146735	Chrysene-d12	21.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p,p'-DDT	352	C14H9Cl5	000050-29-3	83
2			2,2-Bis(p-chlorophenyl)ethanol	266	C14H12Cl2O	002642-82-2	68
3			1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	64
4			m,p'-DDD	318	C14H10Cl4	004329-12-8	62
5			2,2-Bis(p-chlorophenyl)ethanol	266	C14H12Cl2O	002642-82-2	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
Data File : BN019637.D
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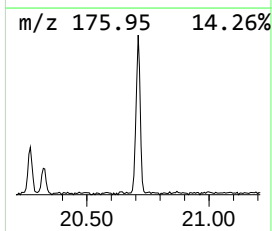
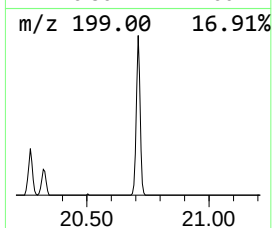
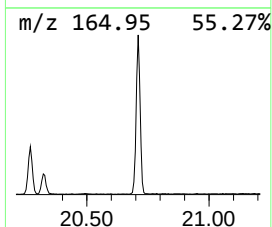
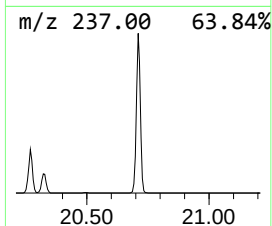
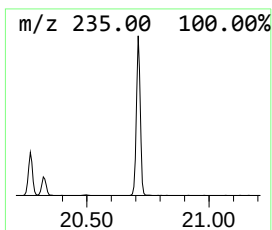
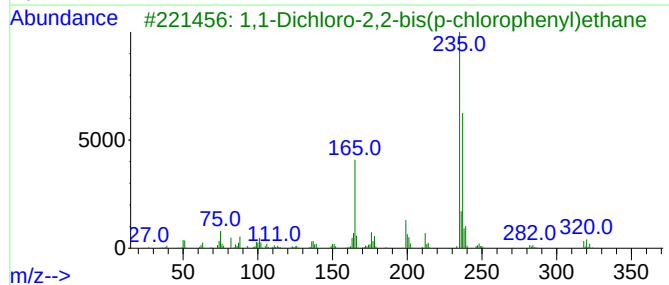
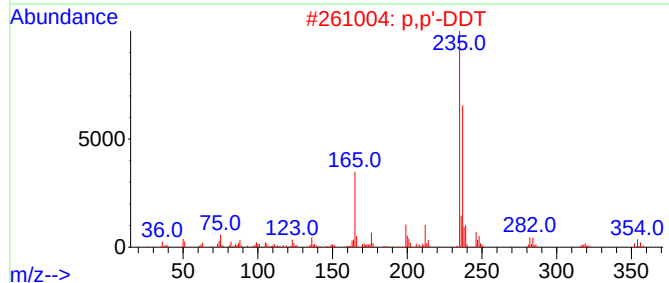
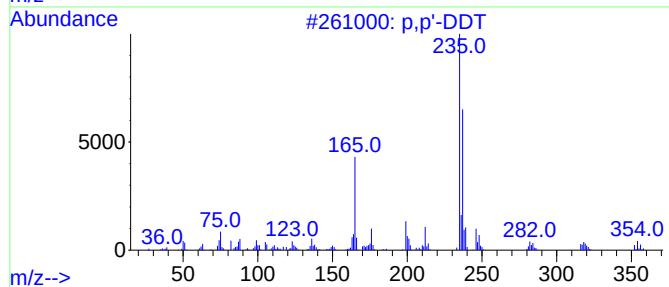
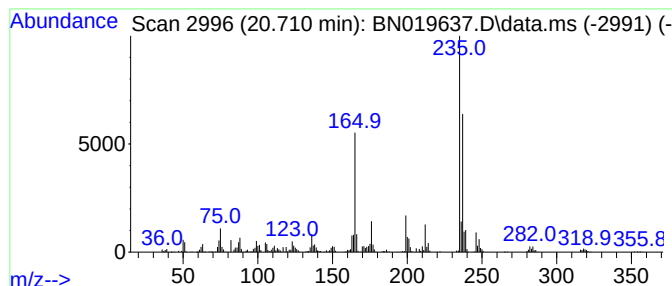
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1,1-Dichloro-2,2-bis(p-chlo... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.710	16.01 ng/ul	983388	Chrysene-d12	21.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p,p'-DDT	352	C14H9Cl5	000050-29-3	99
2			p,p'-DDT	352	C14H9Cl5	000050-29-3	98
3			1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	97
4			Mitotane	318	C14H10Cl4	000053-19-0	97
5			p,p'-DDT	352	C14H9Cl5	000050-29-3	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
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 Acq On : 03 May 2022 01:43
 Operator : CG/JU
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 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ETNS5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	3.281	88.9	ng/u1	4190980	1	7.869	942720	20.0
(DEL) Alkane: C...	3.611	2.9	ng/u1	138356	1	7.869	942720	20.0
1,1'-Biphenyl, ...	17.110	4.1	ng/u1	313844	4	17.227	1544640	20.0
1,1'-Biphenyl, ...	17.404	3.0	ng/u1	227647	4	17.227	1544640	20.0
1,1'-Biphenyl, ...	17.833	6.6	ng/u1	507763	4	17.227	1544640	20.0
1,1'-Biphenyl, ...	19.375	2.5	ng/u1	150358	5	21.410	1228380	20.0
o,p'-DDE	19.763	3.5	ng/u1	213890	5	21.410	1228380	20.0
Mitotane	19.863	2.6	ng/u1	162814	5	21.410	1228380	20.0
m,p'-DDD	20.269	4.6	ng/u1	281078	5	21.410	1228380	20.0
p,p'-DDT	20.322	2.4	ng/u1	146735	5	21.410	1228380	20.0
1,1-Dichloro-2,...	20.710	16.0	ng/u1	983388	5	21.410	1228380	20.0